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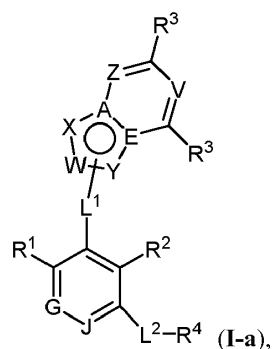
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- (54) Title **NUCLEAR RECEPTOR MODULATORS (ROR) FOR THE TREATMENT OF INFLAMMATORY AND AUTOIMMUNE DISEASES**
- (56) References
Cited: WO-A1-2012/076672
WO-A1-2012/106995
WO-A1-2014/026328
WO-A1-2015/036411
BENJAMIN P. FAUBER ET AL: "Modulators of the Nuclear Receptor Retinoic Acid Receptor-Related Orphan Receptor-[gamma] (ROR[gamma] or RORc)", JOURNAL OF MEDICINAL CHEMISTRY, vol. 57, no. 14, 24 July 2014 (2014-07-24) , pages 5871-5892, XP055242989, US ISSN: 0022-2623, DOI: 10.1021/jm401901d

Enclosed is a translation of the patent claims in Norwegian. Please note that as per the Norwegian Patents Acts, section 66i the patent will receive protection in Norway only as far as there is agreement between the translation and the language of the application/patent granted at the EPO. In matters concerning the validity of the patent, language of the application/patent granted at the EPO will be used as the basis for the decision. The patent documents published by the EPO are available through Espacenet (<http://worldwide.espacenet.com>) or via the search engine on our website here: <https://search.patentstyret.no/>

Patentkrav

1. Forbindelse med formel (I-a):



eller et farmasøytisk akseptabelt salt av samme, hvori:

5 G og J er uavhengig CH eller N;

W er C eller CR^a, L¹ er forbundet med W eller Y; og

A og E er uavhengig C eller N forutsatt at begge ikke er N;

V er CR³ eller N;

X er CR^a, NR^a, N eller O;

10 Y er C, CR^a, NR^a, N, O eller S;

Z er CR³ eller N; eller

W er N eller NR^a, L¹ er forbundet med W eller Y; og

A og E er uavhengig C eller N forutsatt at begge ikke er N;

V er CR³ eller N;

15 X er CR^a, NR^a eller N;

Y er C, CR^a eller N;

Z er CR³ eller N; eller

L¹ er -CH₂- eller -C(O)-;

L² er -C(O)-;

20 R¹ og R² er uavhengig halogen, -O-(C₁-C₃)alkyl, syklopropyl, eller (C₁-C₃)alkyl;

hver R³ er uavhengig H, -CF₃, -CN, halogen, -OCHF₂, -OCF₃, -C(O)N(R^e)₂, -NR^eCOR^e, -(C₁-C₃)alkyl, -O-(C₁-C₃)alkyl, -S-(C₁-C₃)alkyl, (C₃-C₆)sykloalkyl, eller morfolinyl;

- R^4 er (C_1-C_6) alkyl valgfritt substituert med én eller flere substituent uavhengig valgt fra halo, $-CO_2H$, og $-CO_2(C_1-C_4)$ alkyl; eller
- R^4 er $-NH-CH_2$ -tetrahydro-2H-pyranyl valgfritt substituert med én eller to substituent uavhengig valgt fra $-OH$ og $-O(C_1-C_4)$ alkyl; eller
- 5 R^4 er $-N(H)-(C_3-C_6)$ sykloalkyl valgfritt substituert med én eller flere substituent uavhengig valgt fra F, $-OH$, $-COOH$, og $-CO_2(C_1-C_4)$ alkyl; eller
- R^4 er (C_3-C_6) sykloalkyl valgfritt substituert med én eller flere substituent uavhengig valgt fra F, $-OH$, $-COOH$, og $-CO_2(C_1-C_4)$ alkyl; eller
- R^4 er azabisyklo[2,2,1]heptanyl eller azabisyklo[3,1,0]heksanyl, hver valgfritt substituert med $-CO_2H$ eller $-CO_2(C_1-C_4)$ alkyl; eller
- 10 R^4 er azaspiro[3,3]heptanyl valgfritt substituert med $-CH_2CO_2H$, $-CH_2CO_2CH_3$, $-CO_2H$, $-CO_2CH_3$; eller
- R^4 er 2-oxa-6-azaspiro[3,3]heptanyl, 2-oxa-7-azaspiro[3,5]nonanyl, eller 2-oxa-8-azaspiro[4,5]dekanyl, hver valgfritt substituert med $-CH_2CO_2H$; eller
- 15 R^4 er azetidinyll valgfritt substituert med én eller to substituent uavhengig valgt fra $-CH_3$, $-OH$, $-OCH_3$, $-(C_1-C_4)$ alkylen-OH, $-CH_2OCH_3$, $-(C_1-C_4)$ alkylenO (C_1-C_4) alkyl, $-CO_2(C_1-C_4)$ alkyl, $-CO_2H$, $-(C_1-C_4)$ alkylen- CO_2H , $-(C_1-C_4)$ alkylen- $CO_2-(C_1-C_4)$ alkyl, $-N(CH_3)_2$, og pyrrolidinyllgrupper; eller
- R^4 er 1,2-diazepanyl eller 1,4-diazepanyl, hver valgfritt substituert med $-(C_1-C_4)$ alkylen-OH; eller
- 20 R^4 er morfolinyll valgfritt substituert med $=O$; eller
- R^4 er piperazinyll, piperidinyll, eller pyrrolidinyll, hver valgfritt substituert med én eller flere substituent uavhengig valgt fra halo, $-CN$, $-(C_1-C_4)$ alkyl, $-CH_2$ -syklopropyl, $-(C_1-C_4)$ alkylen-F, $-CF_3$, $-CH_2CF_3$, $-COOH$, $-(C_1-C_4)$ alkylenCOOH, $-CH(OH)CO_2H$, $-COCH_3$, $-CO_2(C_1-C_4)$ alkyl, $-CH(OH)CO_2CH_3$, $-(C_1-C_4)$ alkylen-C(=O)O (C_1-C_4) alkyl, $-(C_1-C_4)$ alkylen-OH, $-OH$, $-O(C_1-C_4)$ alkyl, $-(C_1-C_4)$ alkylenO (C_1-C_4) alkyl, $-O(C_1-C_4)$ alkylenCO $_2H$, $-O(C_1-C_4)$ alkylenC(=O)O (C_1-C_4) alkyl, $-CONHCH_3$, $-SO_2-(C_1-C_4)$ alkyl, $-NH_2$, $-NH(C_1-C_4)$ alkyl, $-N((C_1-C_4)alkyl)_2$, syklobutankarboksylsyre, og oksetanyl;
- 25 hver R^a er uavhengig H, $-C(O)CH_3$, valgfritt substituert (C_1-C_6) alkyl, valgfritt substituert (C_3-C_6) sykloalkyl eller $-S(O)_2$ -fenyl; og
- 30 R^e er uavhengig H eller $-(C_1-C_3)$ alkyl;

forutsatt at ikke flere enn to av A, E, W, X og Y er N.

2. Forbindelse ifølge krav 1, eller et farmasøytisk akseptabelt salt av samme, hvori R^1 og R^2 begge er halo eller $-CH_3$.
3. Forbindelse ifølge ett av kravene 1 eller 2, eller et farmasøytisk akseptabelt salt av samme, hvori
5 W er C eller CH, A er C, og E er C.
4. Forbindelse ifølge ett av kravene 1 til 3, eller et farmasøytisk akseptabelt salt av samme, hvori G er CH.
5. Forbindelse ifølge ett av kravene 1 til 4, eller et farmasøytisk akseptabelt salt av samme, hvori J er CH.
- 10 6. Forbindelse ifølge krav 1, eller et farmasøytisk akseptabelt salt av samme, hvori:
 - W er CH;
 - X er CR^a , hvori R^a er H eller $-CH_3$;
 - Y er N;
 - A er C;
 - 15 E er C;
 - V er CR^3 , hvori R^3 er H;
 - Z er N eller CR^3 , hvori R^3 er H, $-CH_3$, eller $-CF_3$;
 - L^1 er forbundet med Y;
 - L^1 er $-CH_2-$ eller $-C(O)-$;
 - 20 L^2 er $-C(O)-$;
 - R^1 er $-Cl$ eller $-CH_3$;
 - R^2 er $-Cl$ eller $-CH_3$;
 - hver R^3 er uavhengig H, $-F$, $-Cl$, $-CN$, $-CH_3$, $-CH(CH_3)_2$, $-CF_3$, $-OCHF_2$, $-OCF_3$, $-SCH_3$, eller syklopropyl; og
 - 25 R^4 er valgfritt substituert 2-oxa-6-azaspiro[3,3]heptanyl eller valgfritt substituert piperidinyll.
7. Forbindelse ifølge krav 6, eller et farmasøytisk akseptabelt salt av samme, hvori det valgfritt substituerte piperidinyll er valgfritt substituert med $-CH_3$, $-COOH$, $-CH_2CO_2H$, $-CH_2CH_2CO_2H$, $-CO_2CH_3$, $-CO_2CH_2CH_3$, eller $-OH$.
8. Forbindelse ifølge krav 1, eller et farmasøytisk akseptabelt salt av samme, hvori:
 - 30 W er C;
 - X er NR^a , hvori R^a er H, $-CH_3$, $-CH_2CH_3$, $-COCH_3$, eller $-CH_2CH(CH_3)_2$;
 - Y er CR^a , hvori R^a er H;

- Z er CR^3 , hvori R^3 er H eller $-CH_3$;
- A er C;
- E er C;
- V er CR^3 , hvori R^3 er H, $-CN$, $-CH_3$, eller $-CF_3$;
- 5 L^1 er forbundet med W;
- L^1 er $-CH_2-$ eller $-C(O)-$;
- L^2 er $-C(O)-$;
- R^1 er $-Cl$, $-CH_3$, eller syklopropyl;
- R^2 er $-Cl$ eller $-CH_3$;
- 10 hver R^3 er uavhengig H, $-F$, $-Br$, $-Cl$, $-CN$, $-CH_3$, $-CF_3$, $-OCH_3$, $-OCHF_2$, $-OCF_3$, $-CONHCH_3$, $-CON(CH_3)_2$, $-NHCOCH_3$, syklopropyl, eller morfolinyl;
- R^4 er valgfritt substituert (C_1-C_6)alkyl, $-NH-CH_2-$ valgfritt substituert tetrahydro-2H-pyranyl, $-NH-$ valgfritt substituert sykloheksyl, valgfritt substituert azabisyklo[2,2,1]heptanyl, valgfritt substituert 3-azabisyklo[3,1,0]heksanyl, valgfritt substituert azaspiro[3,3]heptanyl, valgfritt substituert azetidiny, valgfritt substituert syklobutyl, valgfritt substituert sykloheksyl,
- 15 valgfritt substituert syklopentyl, valgfritt substituert 1,4-diazepanyl, valgfritt substituert morfolinyl, valgfritt substituert 2-oxa-6-azaspiro[3,3]heptanyl, valgfritt substituert 2-oxa-7-azaspiro[3,5]nonanyl, valgfritt substituert 2-oxa-8-azaspiro[4,5]dekanyl, valgfritt substituert piperaziny, valgfritt substituert piperidiny, eller valgfritt substituert pyrrolidiny.
- 20 9. Forbindelse ifølge krav 1, eller et farmasøytisk akseptabelt salt av samme, hvori:
- W er C;
- X er NR^a , hvori R^a er H eller $-CH_3$;
- Y er CR^a , hvori R^a er H;
- Z er N;
- 25 A er C;
- E er C;
- V er CR^3 , hvori R^3 er H;
- L^1 er forbundet med W;
- L^1 er $-CH_2-$ eller $-C(O)-$;
- 30 L^2 er $-C(O)-$;
- R^1 er $-Cl$;
- R^2 er $-Cl$;
- hver R^3 er uavhengig H, $-CH_3$, eller $-CF_3$; og

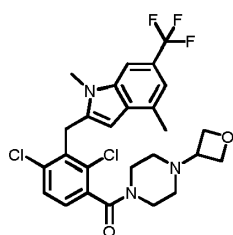
R^4 er valgfritt substituert morfolinyl, valgfritt substituert piperazinyl, eller valgfritt substituert piperidinyl.

10. Forbindelse ifølge krav 1, eller et farmasøytisk akseptabelt salt av samme, hvori:

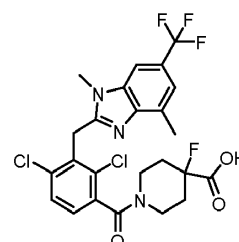
- 5 W er CR^a , hvori R^a er H;
 X er O eller NR^a , hvori R^a er H eller $-CH_3$;
 Y er C;
 Z er CR^3 , hvori R^3 er H eller $-CF_3$;
 A er C;
 E er C;
 10 V er CR^3 , hvori R^3 er H;
 L^1 er forbundet med Y;
 L^1 er $-CH_2-$;
 L^2 er $-C(O)-$;
 R^1 er $-Cl$;
 15 R^2 er $-Cl$;
 hver R^3 er uavhengig H, $-CH_3$, $-CF_3$, eller $-OCF_3$; og
 R^4 er valgfritt substituert piperidinyl.

11. Forbindelse ifølge krav 10 eller et farmasøytisk akseptabelt salt av samme, hvori det valgfritt substituerte piperidinyl er valgfritt substituert med $-CH_3$, $-COOH$, $-CH_2COOH$, $-CH_2CO_2CH_3$, eller $-CO_2CH_2CH_3$.
- 20

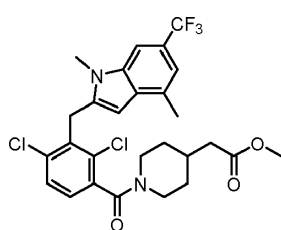
12. Forbindelse ifølge krav 1, hvori forbindelsen er:



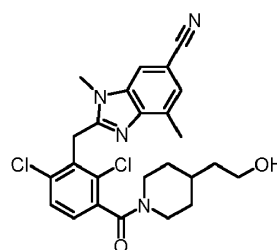
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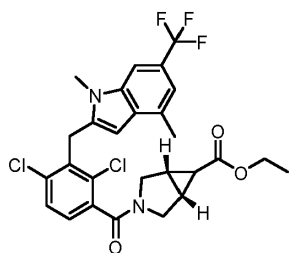
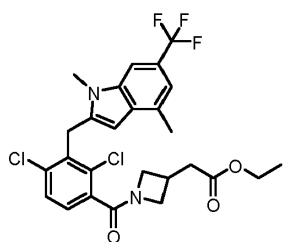
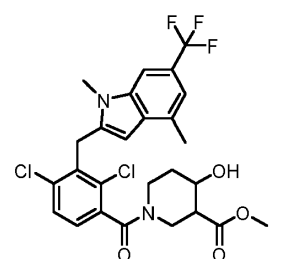
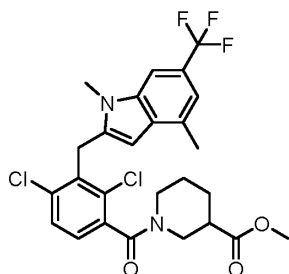
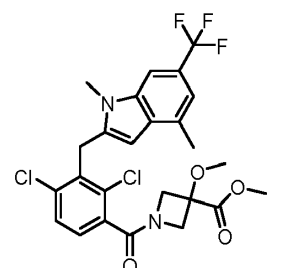
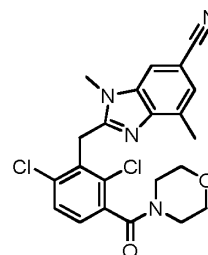
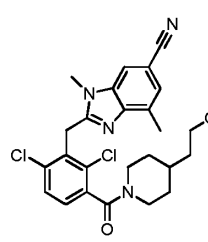
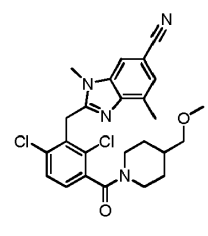
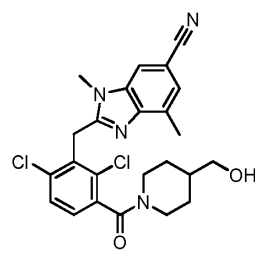
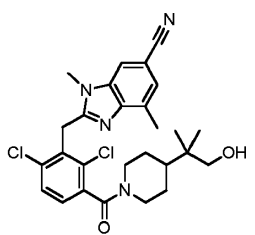
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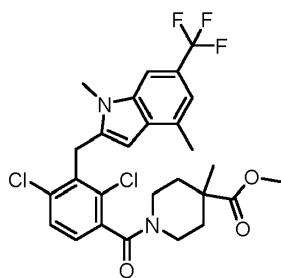
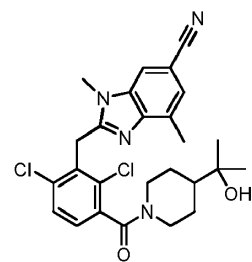
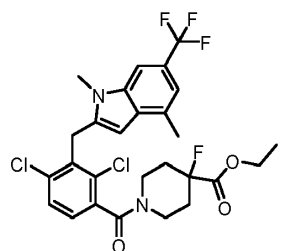
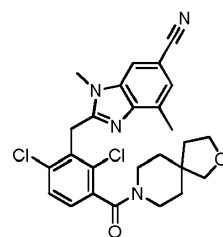
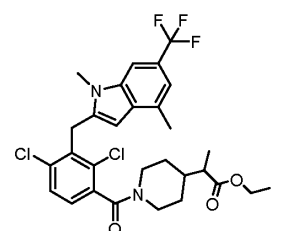
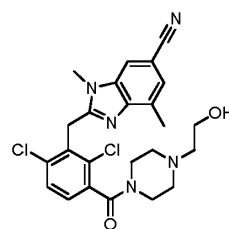
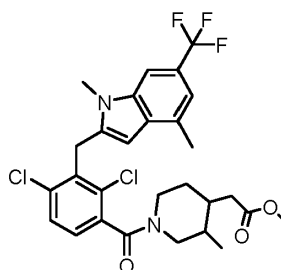
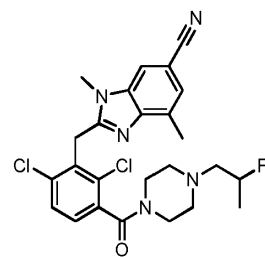
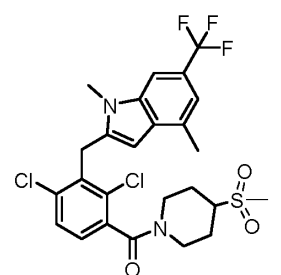
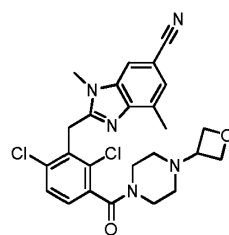


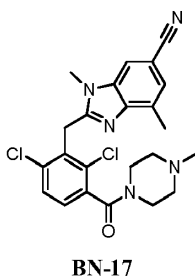
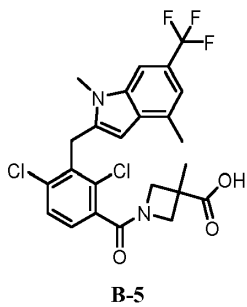
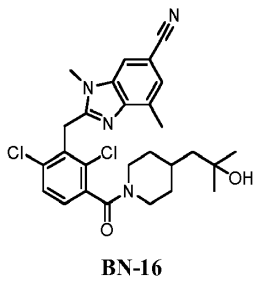
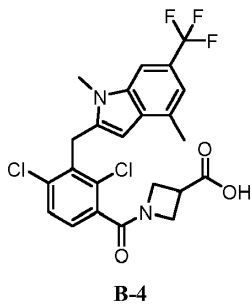
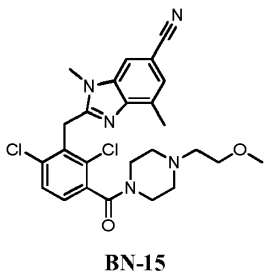
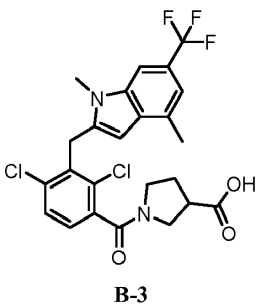
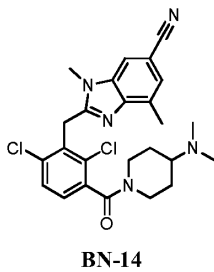
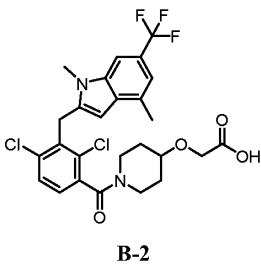
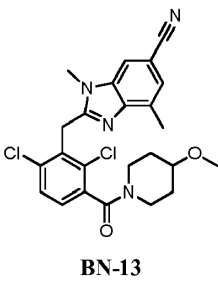
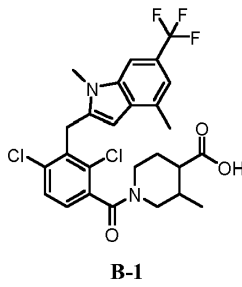
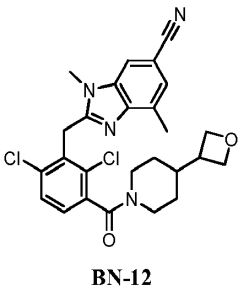
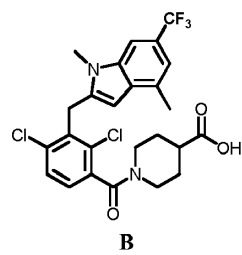
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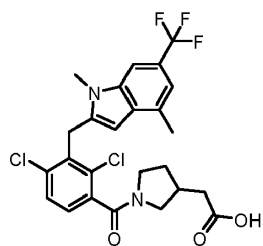
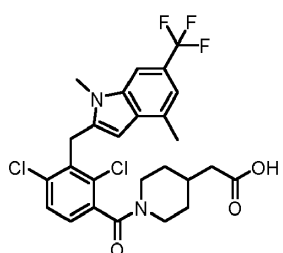
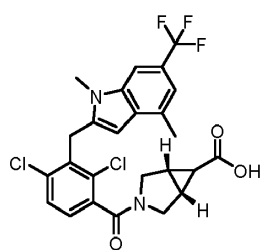
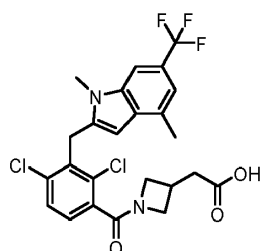
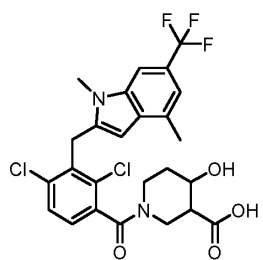
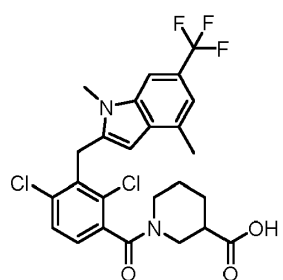
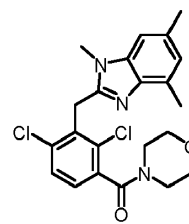
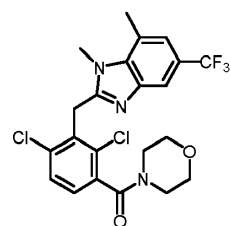
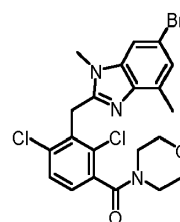
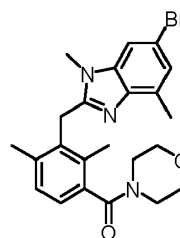
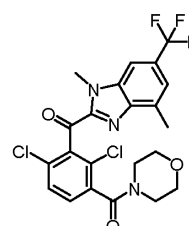
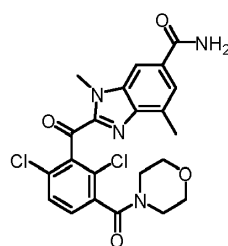


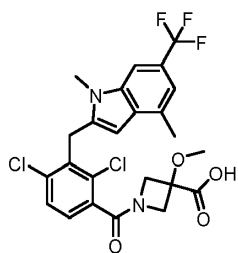
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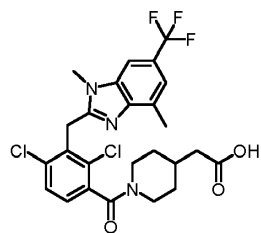
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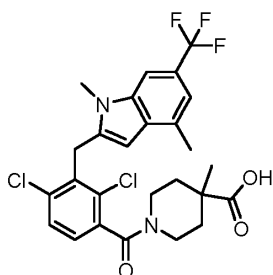
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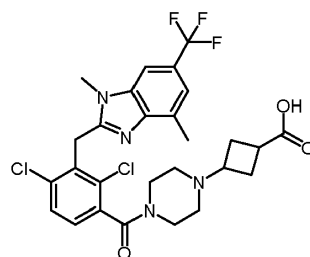
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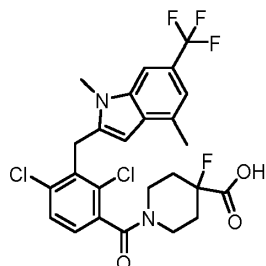
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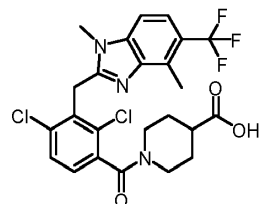
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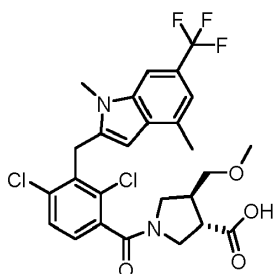
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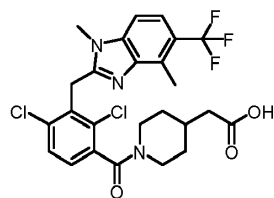
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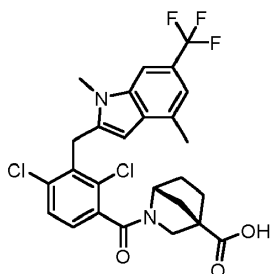
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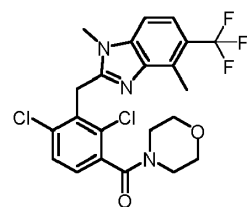
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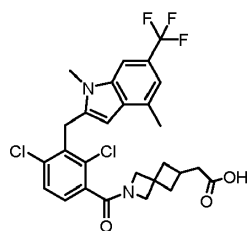
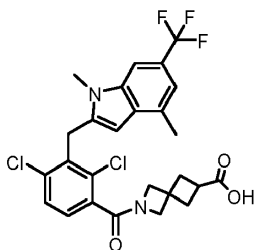
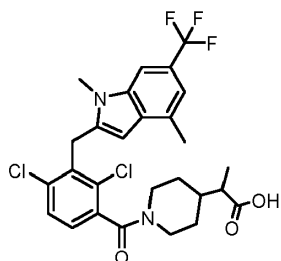
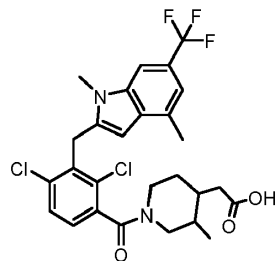
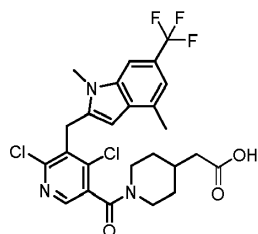
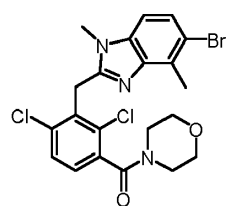
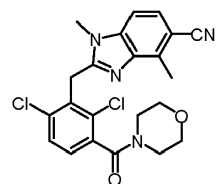
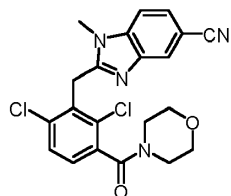
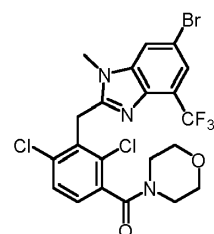
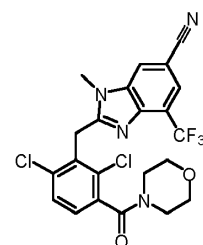
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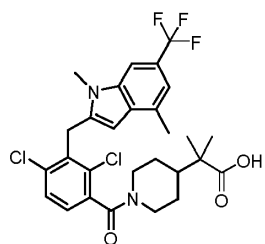
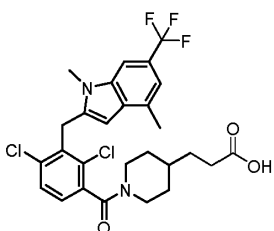
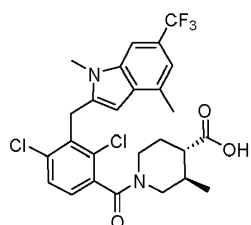
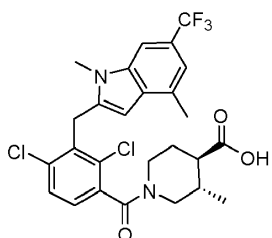
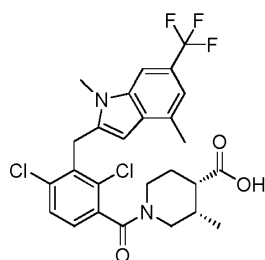
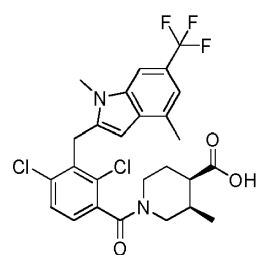
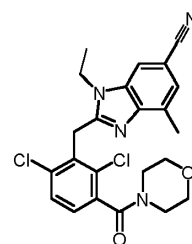
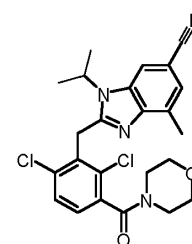
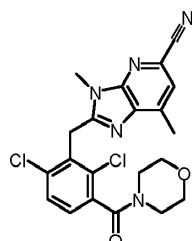
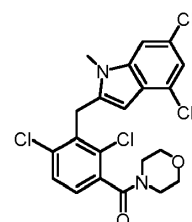
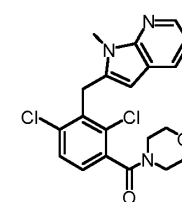
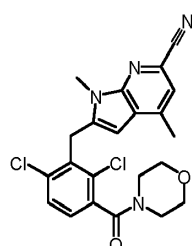
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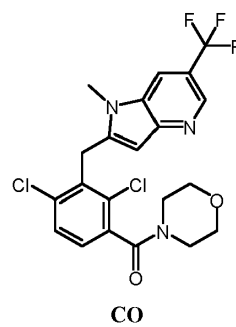
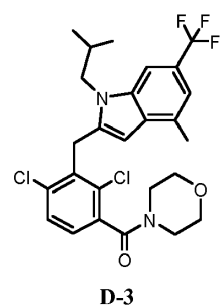
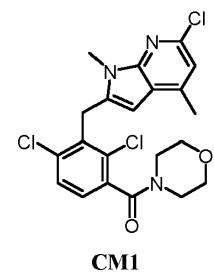
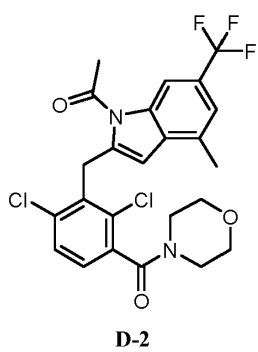
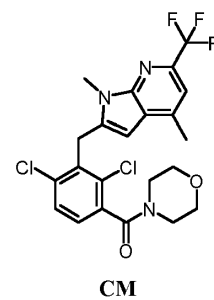
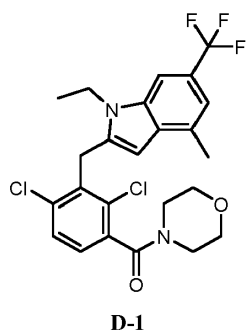
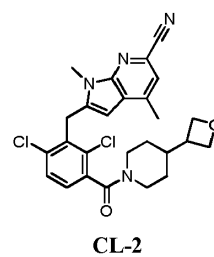
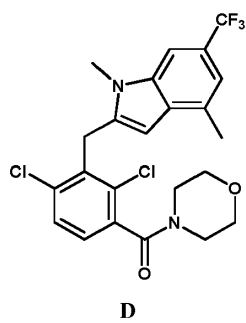
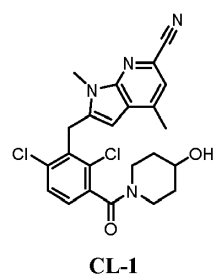
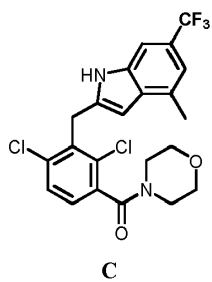
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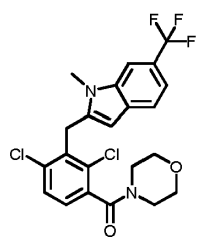
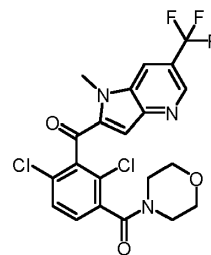
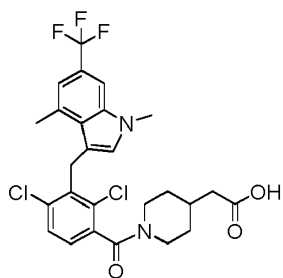
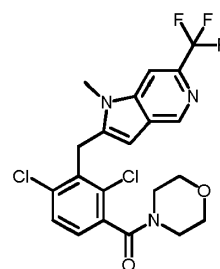
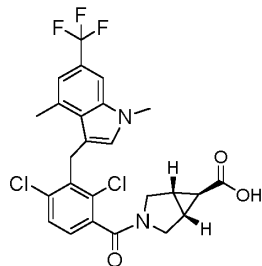
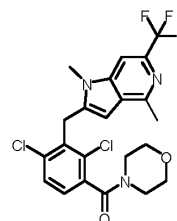
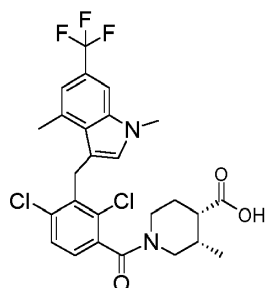
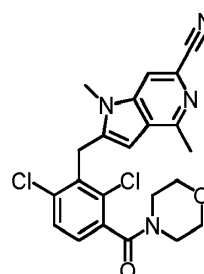
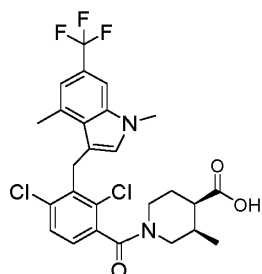
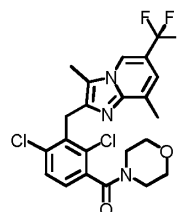
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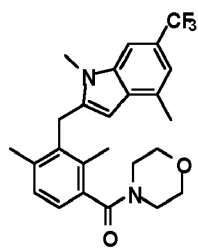
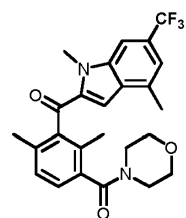
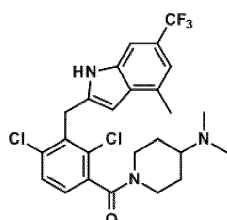
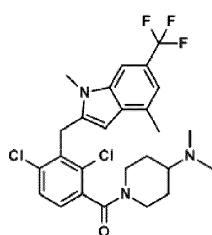
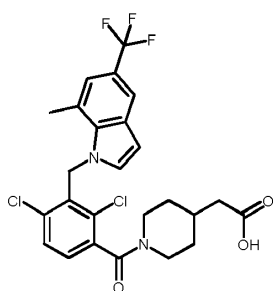
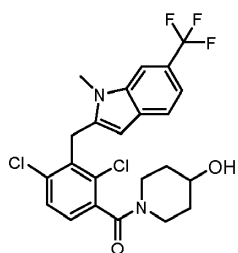
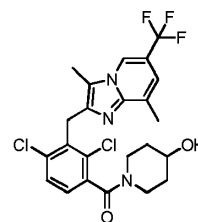
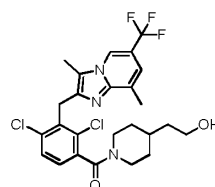
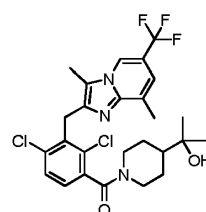
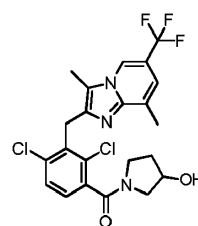
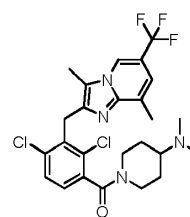
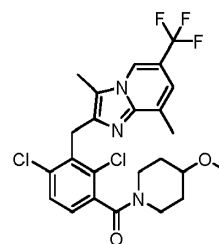
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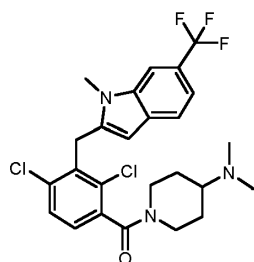
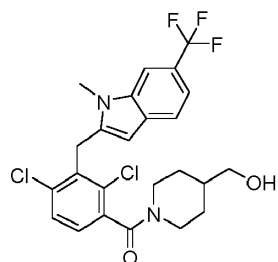
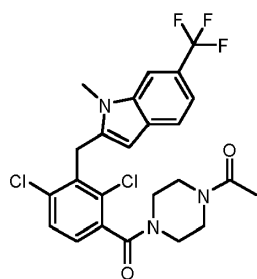
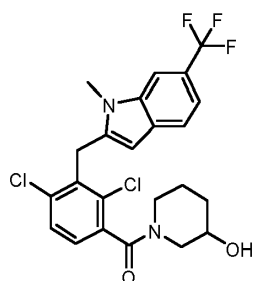
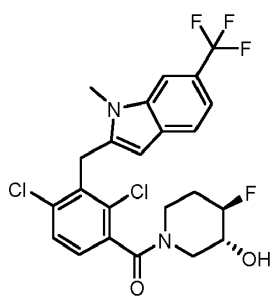
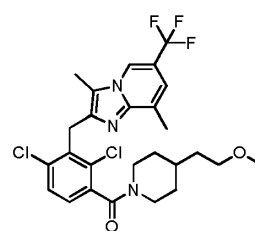
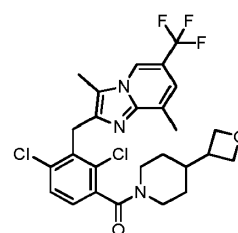
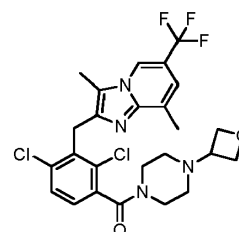
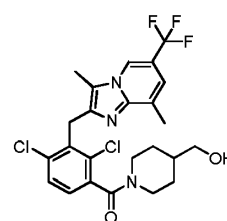
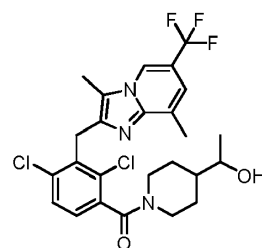
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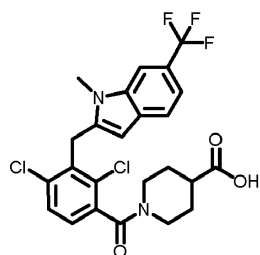
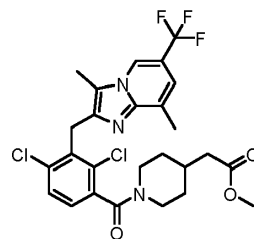
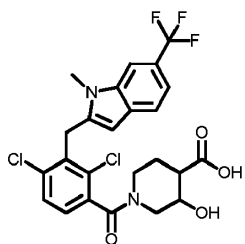
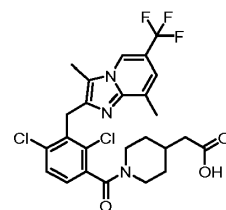
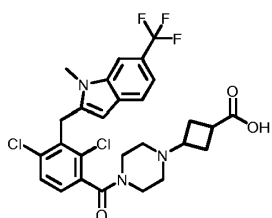
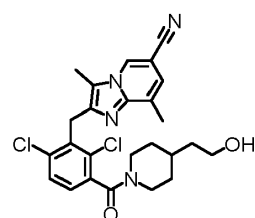
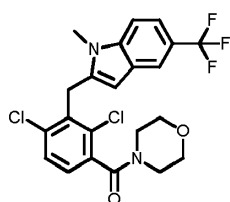
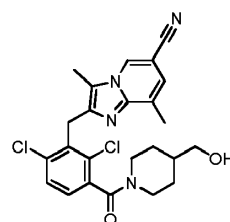
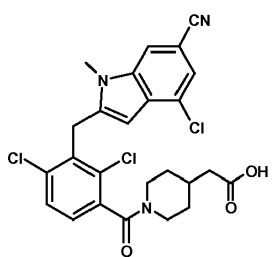
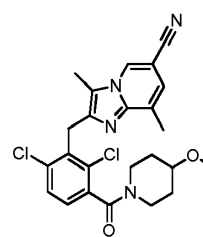
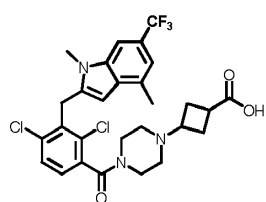
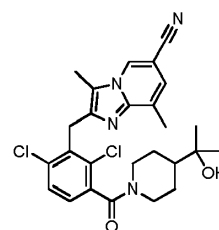
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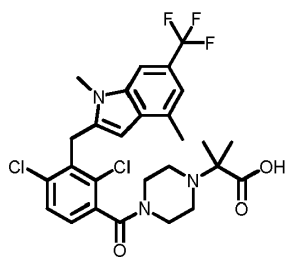
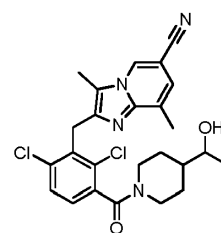
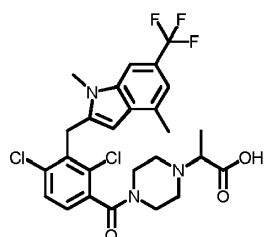
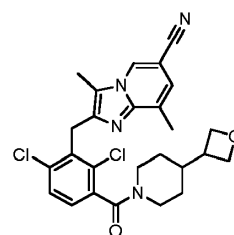
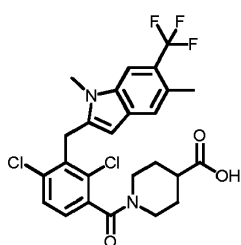
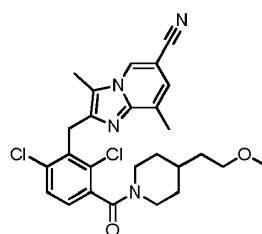
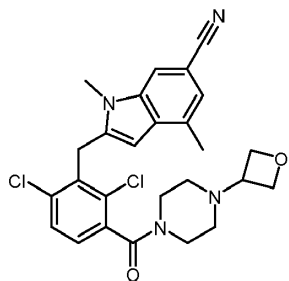
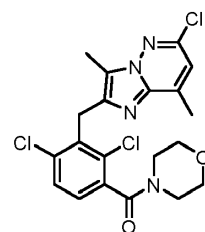
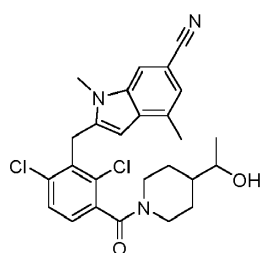
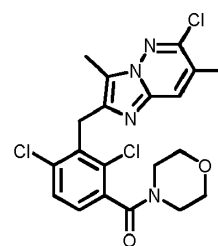
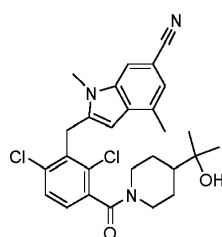
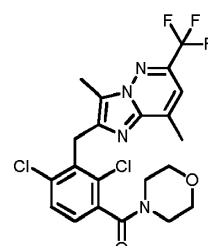


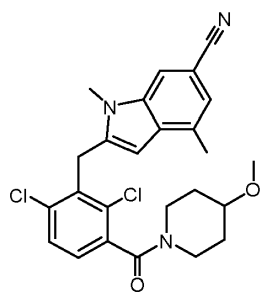
**E****CP****F****CQ****F1-1****CR****F1-2****CS****F1-3****CT**

**G****G1****H****I****J****L****CT-1****CT-2****CT-3****CT-4****CT-5****CT-6**

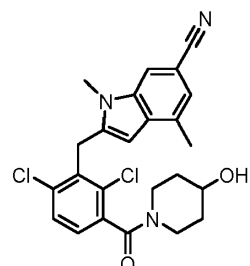
**L-1****L-2****L-3****L-4****L-5****CT-7****CT-8****CT-9****CT-10****CT-11**

**M****CT-12****N****CU****O****CV****P****CV-1****R****CV-2****S****CV-3**

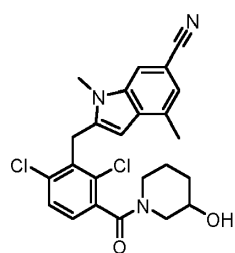
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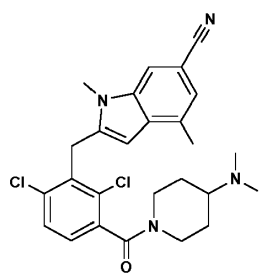
AA-3



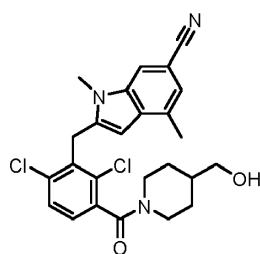
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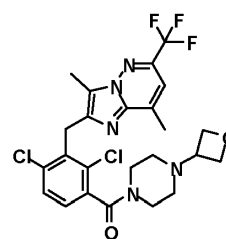
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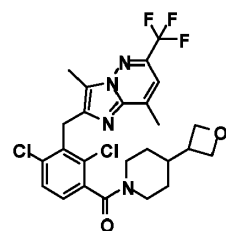
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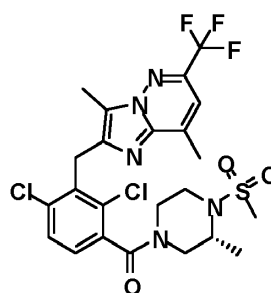
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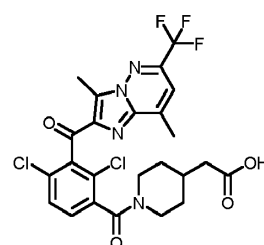
CZ-1



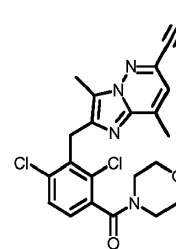
CZ-2



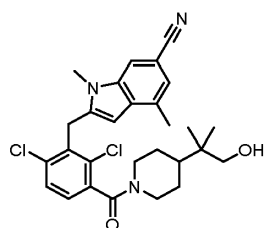
CZ-3



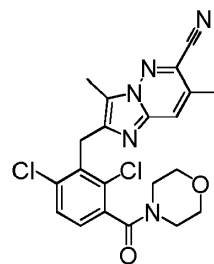
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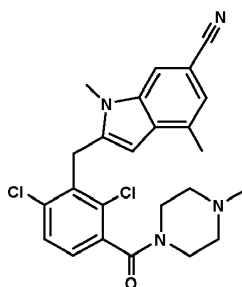
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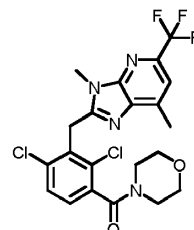
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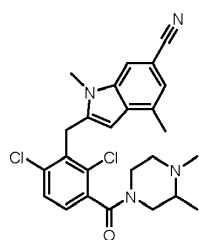
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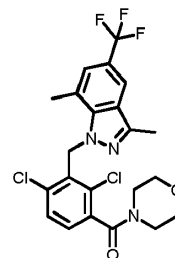
AA-9



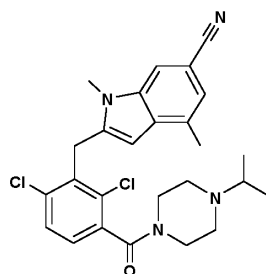
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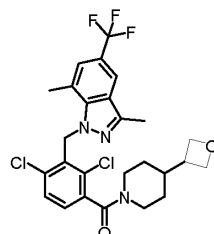
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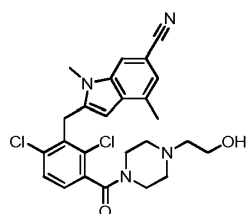
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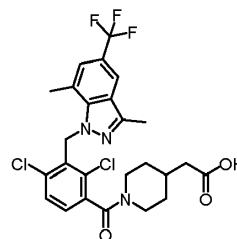
AA-11



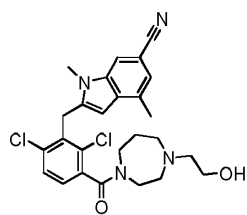
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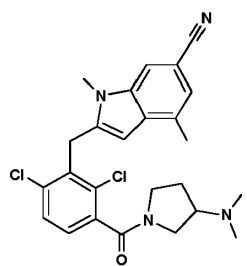
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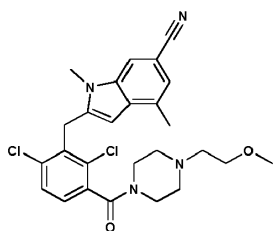
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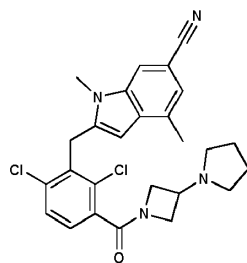
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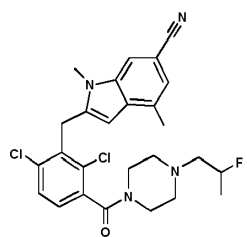
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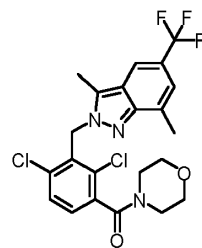
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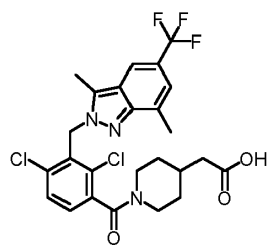
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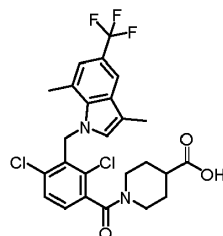
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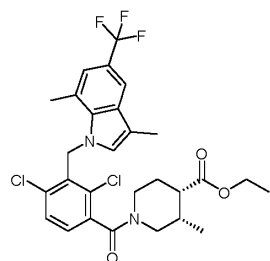
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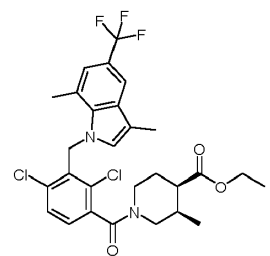
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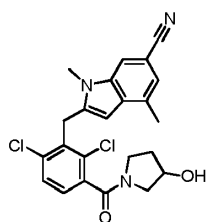
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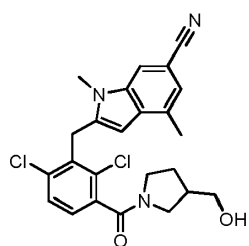
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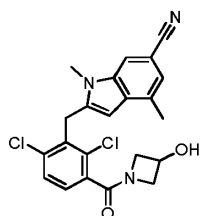
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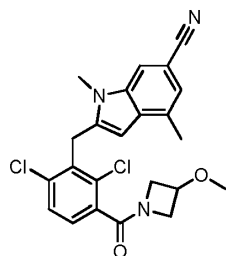
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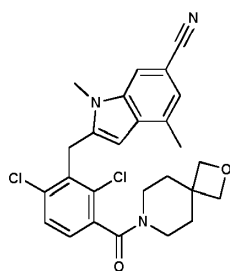
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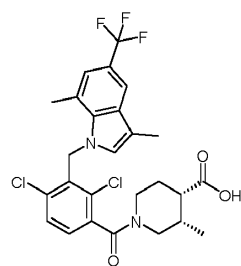
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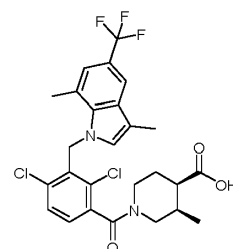
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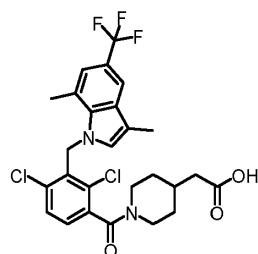
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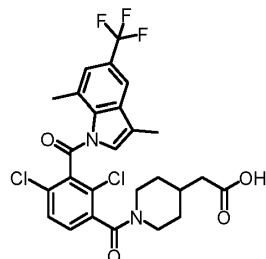
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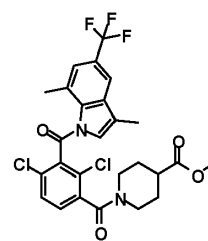
DH1-2



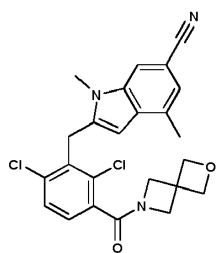
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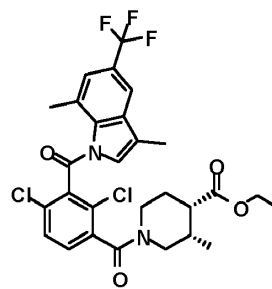
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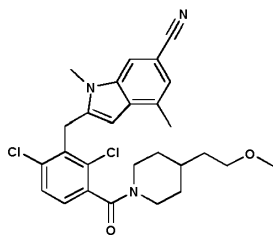
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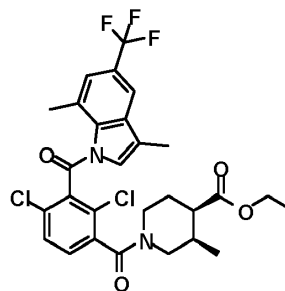
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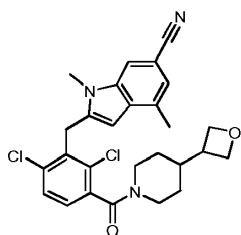
DJ-2



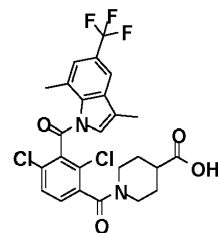
AA-27



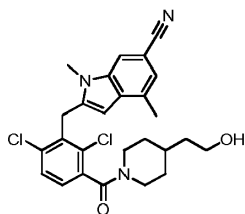
DJ-3



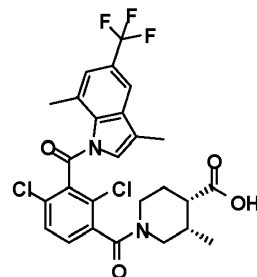
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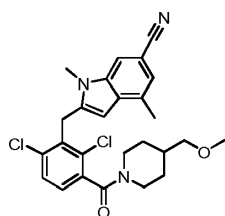
DJ1-1



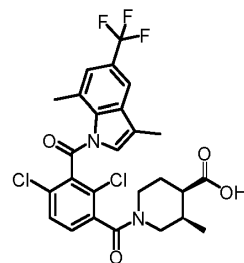
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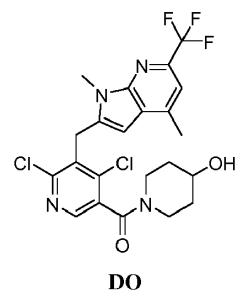
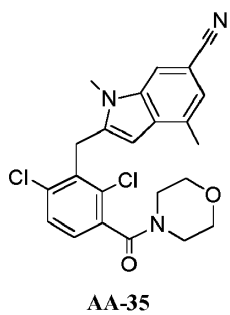
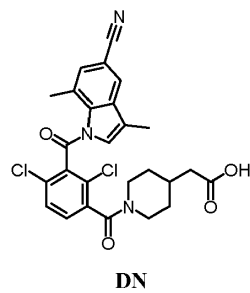
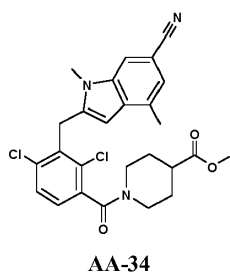
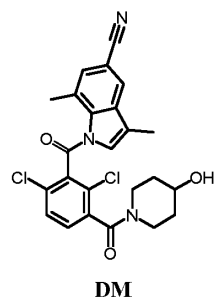
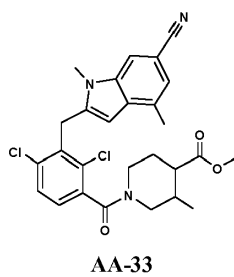
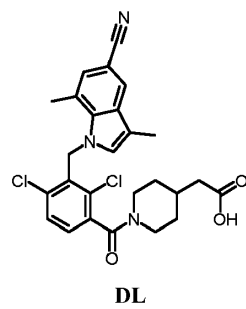
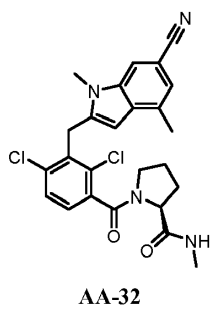
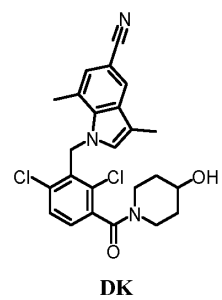
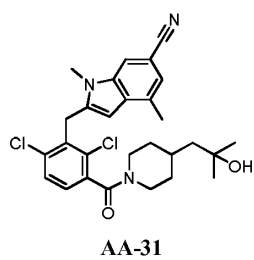
DJ1-2

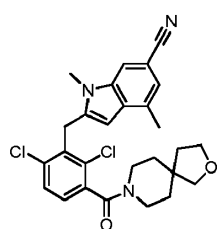


AA-30

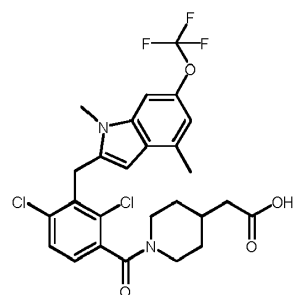


DJ1-3

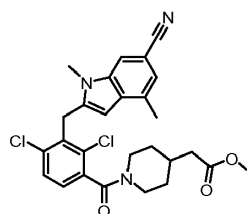




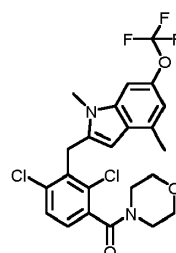
AA-36



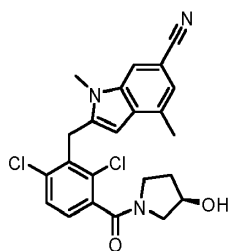
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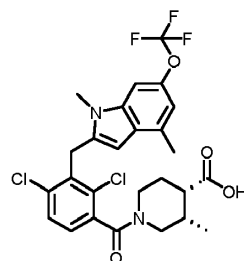
AA-38



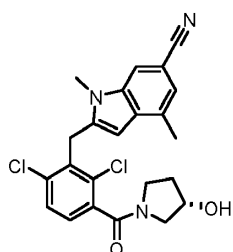
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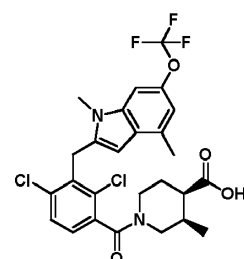
AA-39



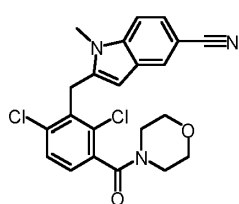
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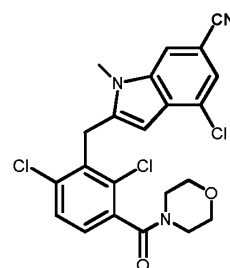
AA-40



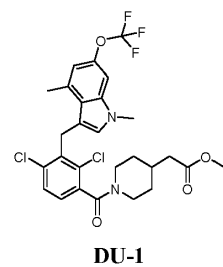
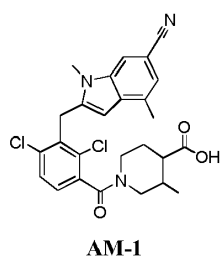
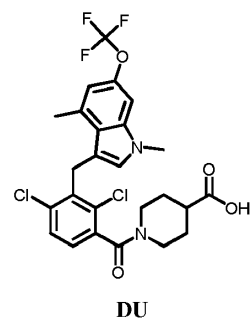
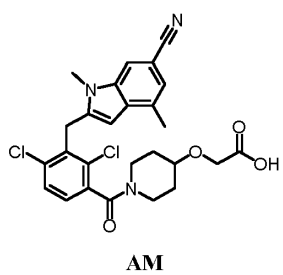
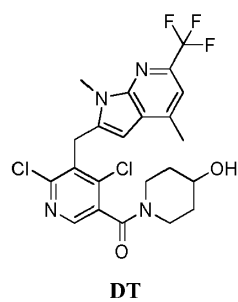
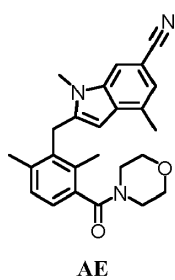
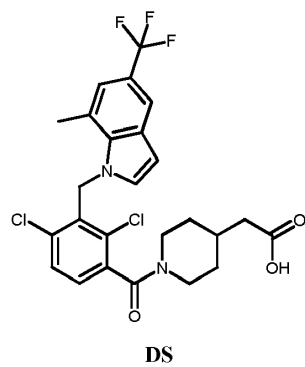
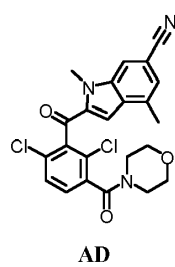
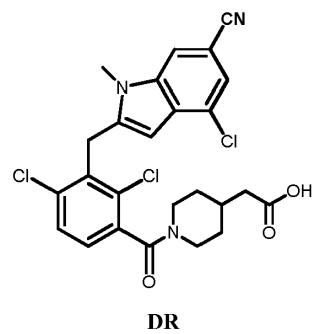
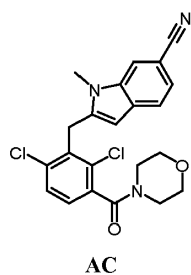
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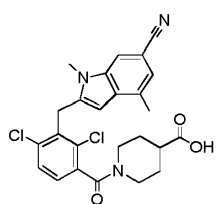
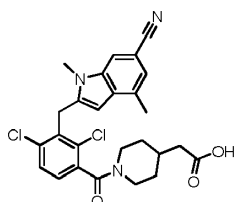
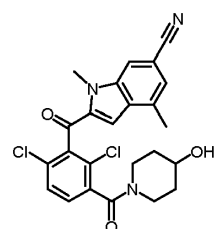
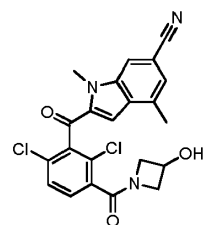
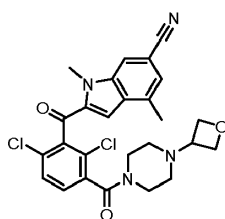
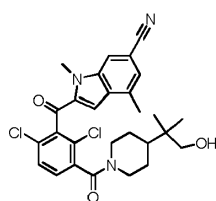
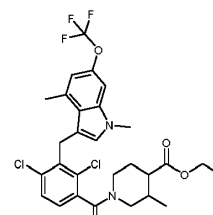
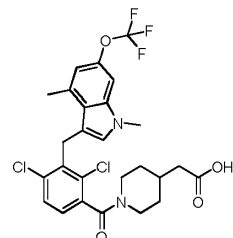
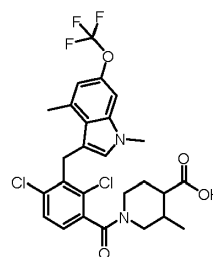
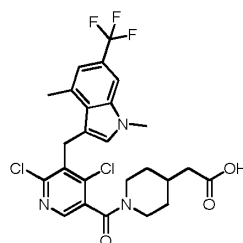
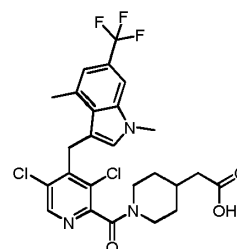
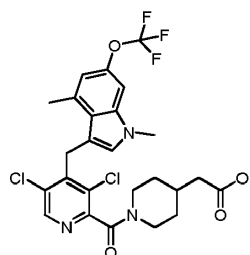


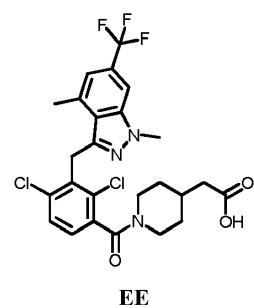
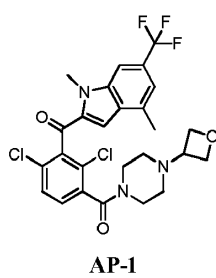
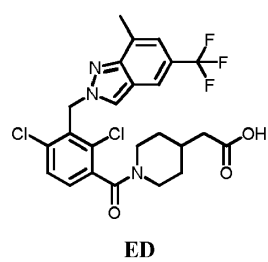
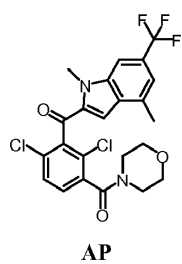
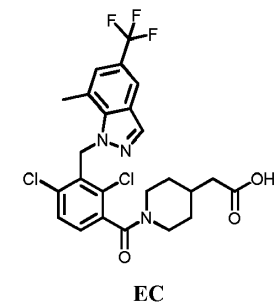
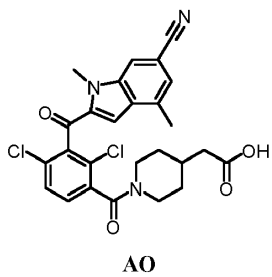
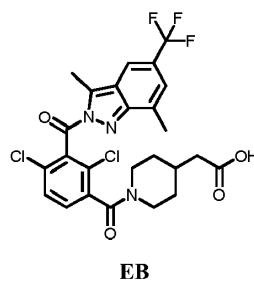
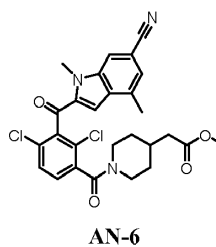
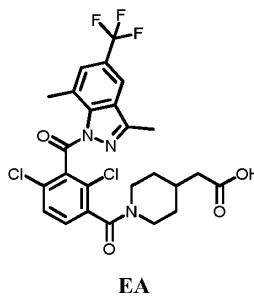
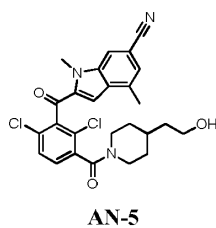
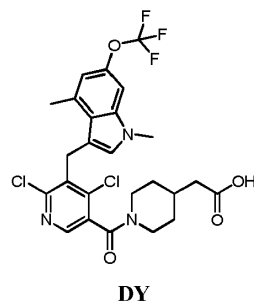
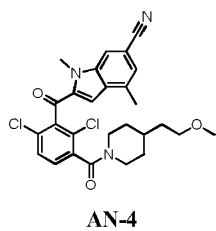
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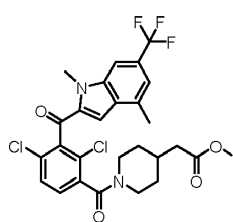
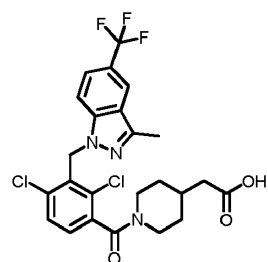
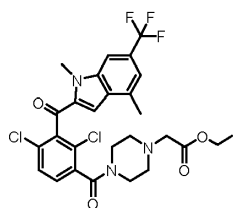
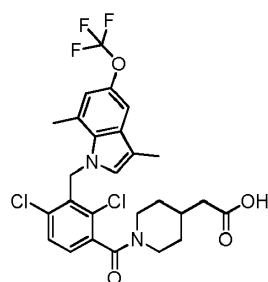
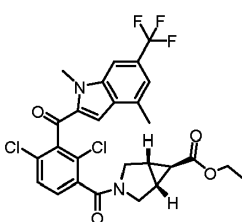
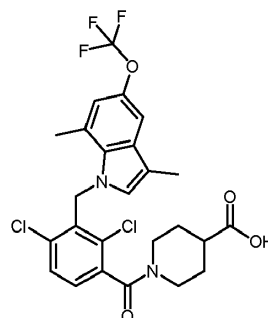
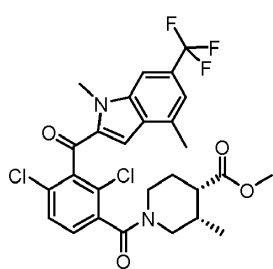
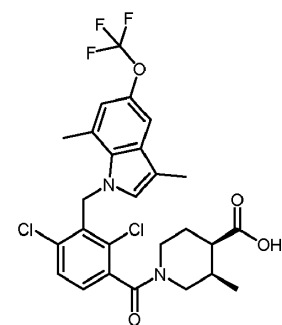
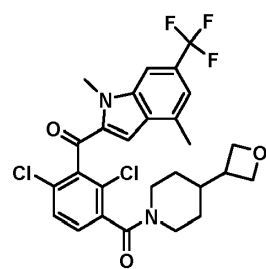
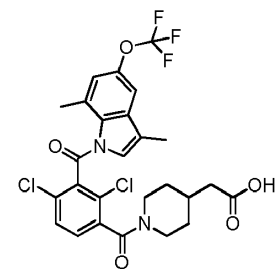


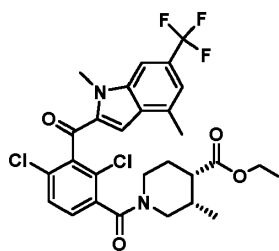
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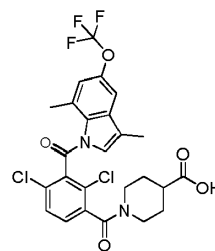
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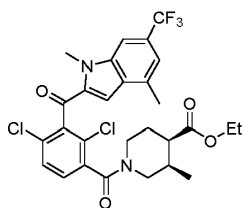
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**AP-7**

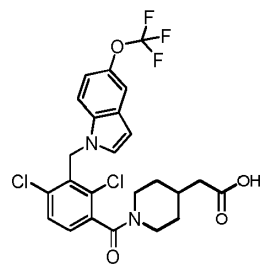
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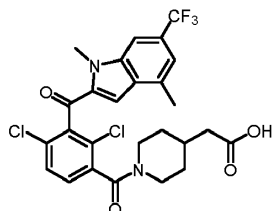
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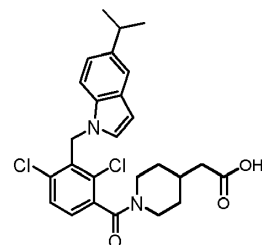
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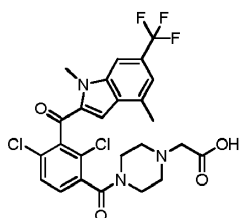
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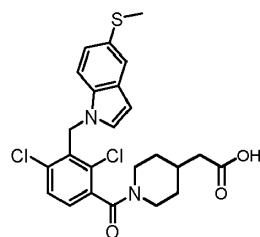
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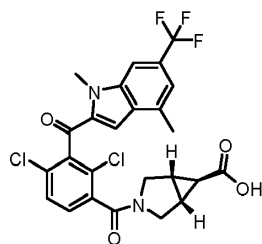
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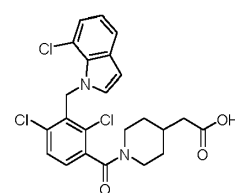
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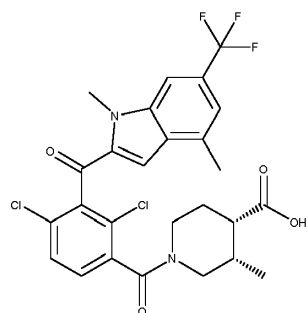
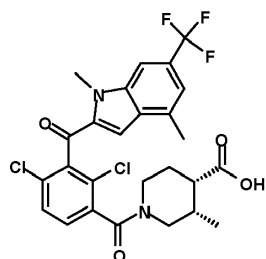
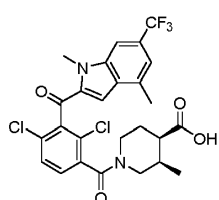
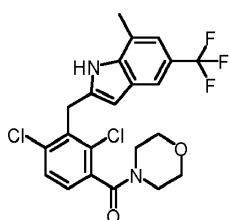
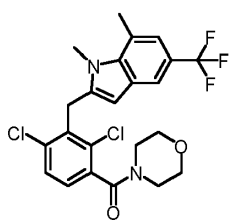
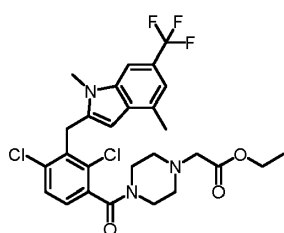
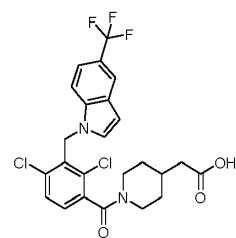
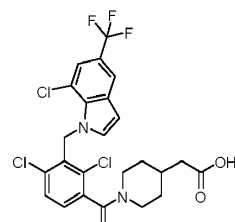
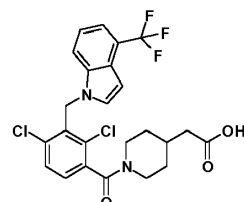
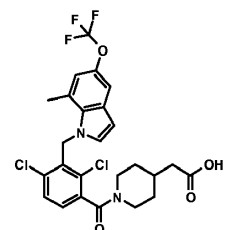
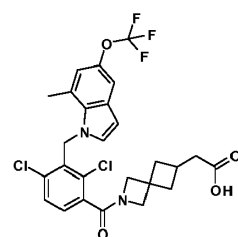
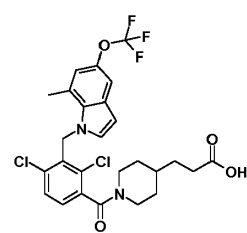
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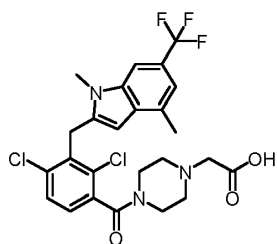
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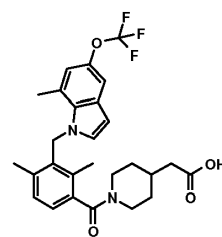
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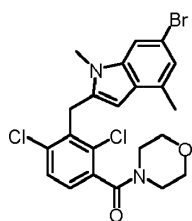
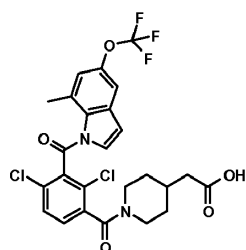
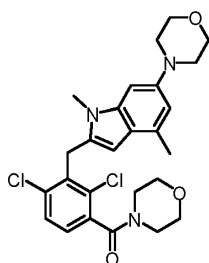
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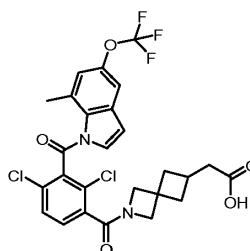
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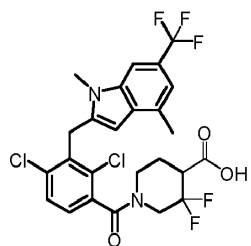
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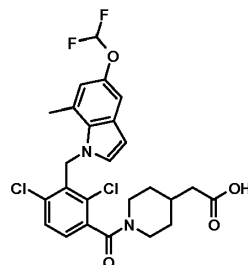
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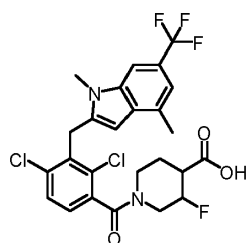
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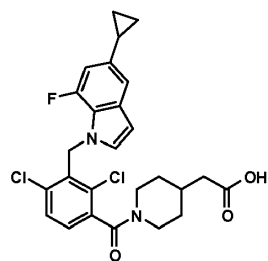
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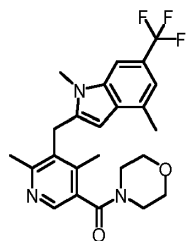
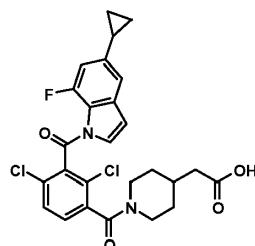
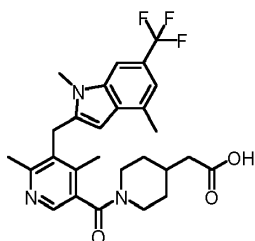
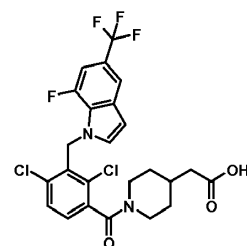
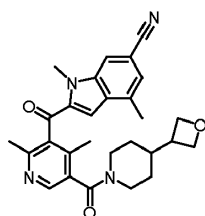
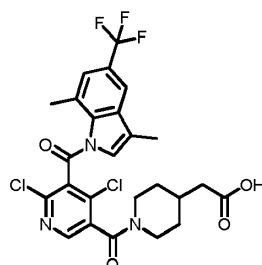
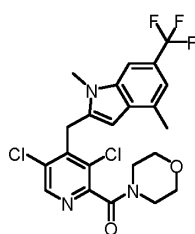
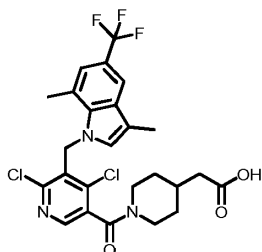
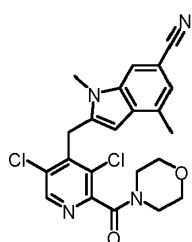
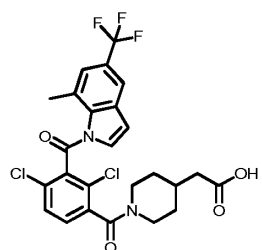
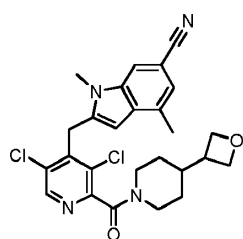
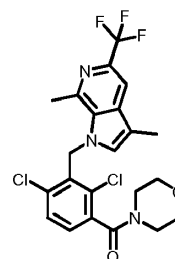
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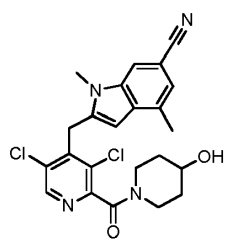
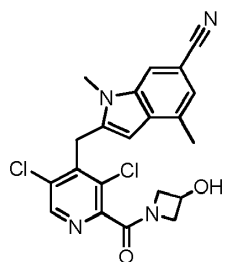
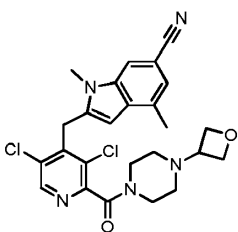
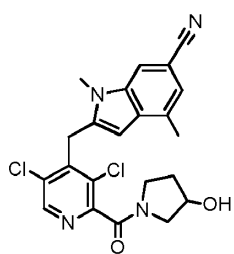
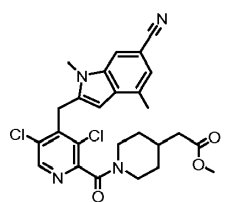
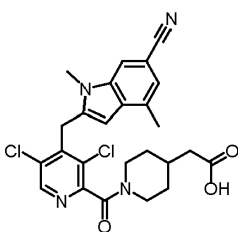
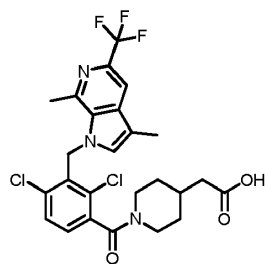
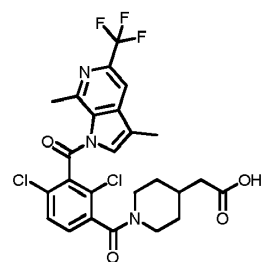
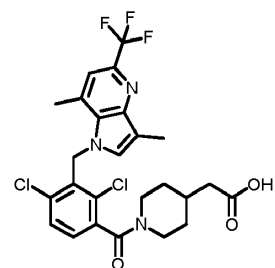
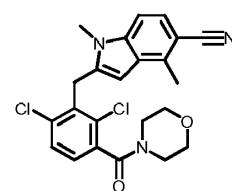
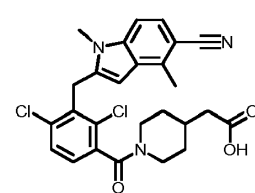
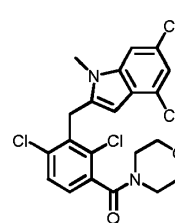


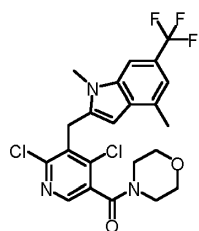
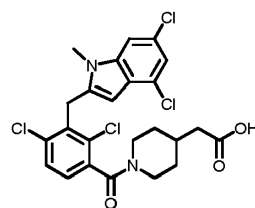
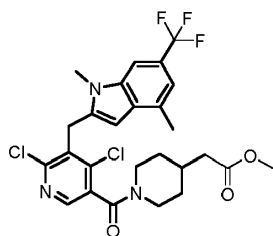
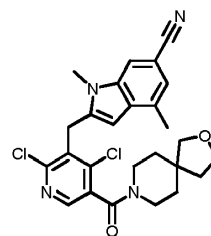
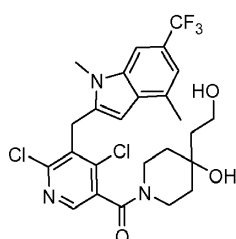
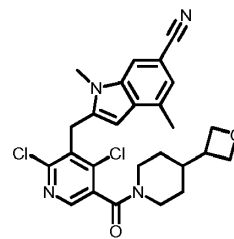
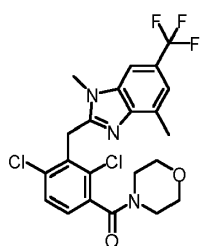
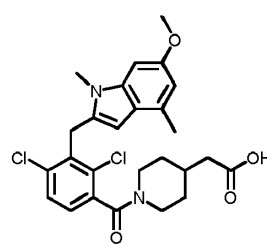
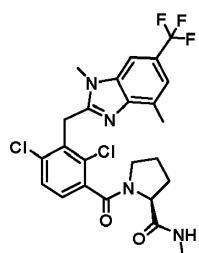
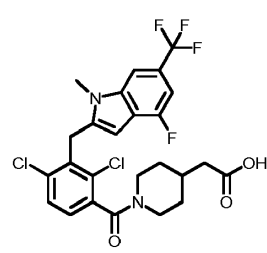
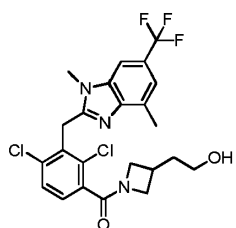
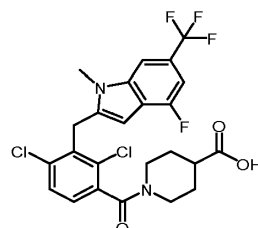
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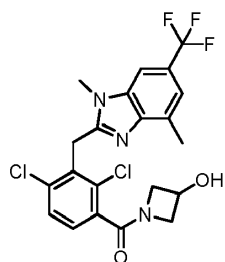
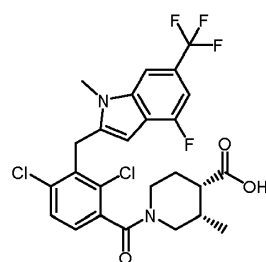
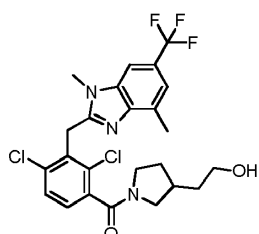
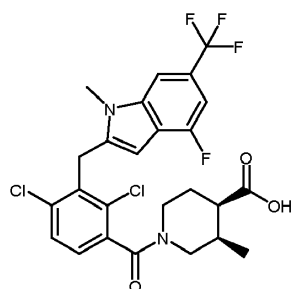
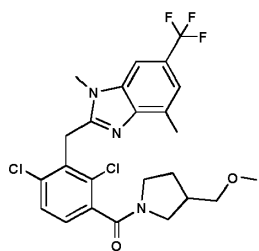
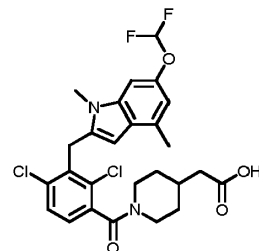
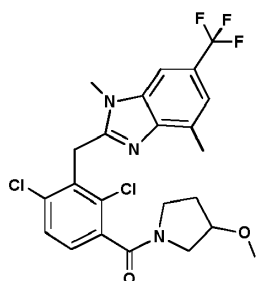
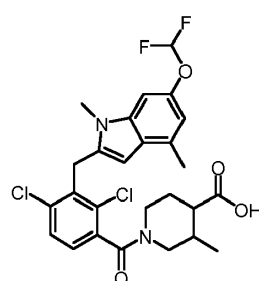
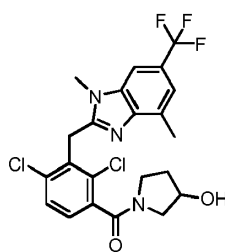
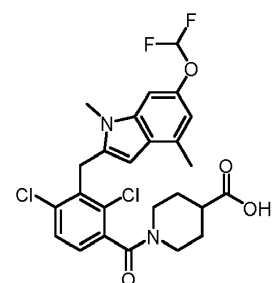


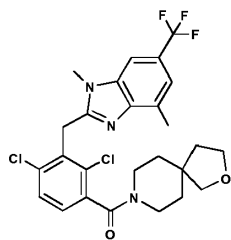
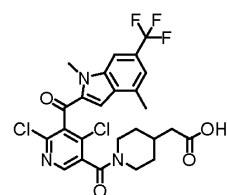
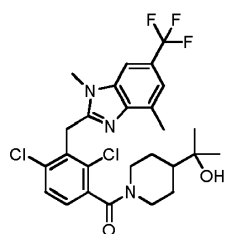
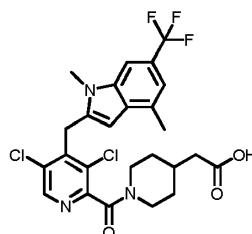
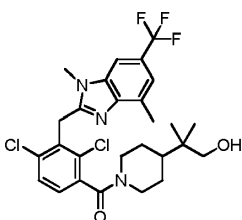
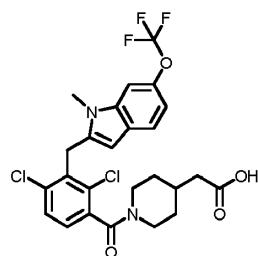
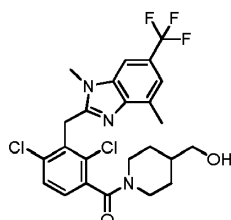
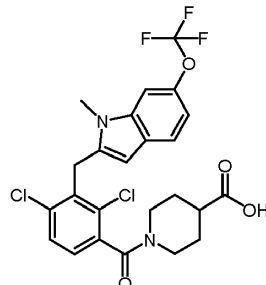
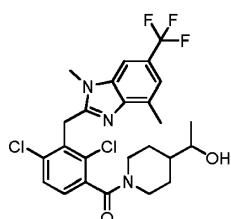
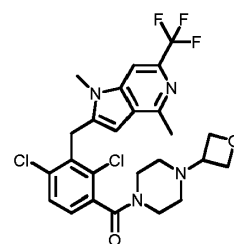
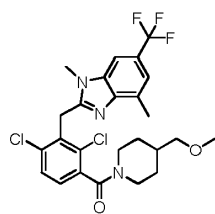
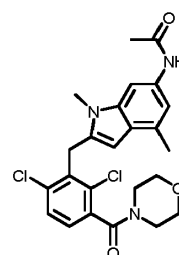
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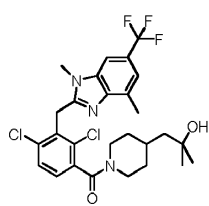
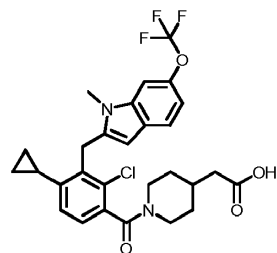
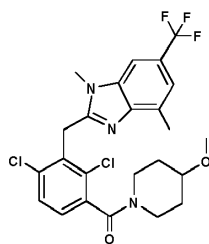
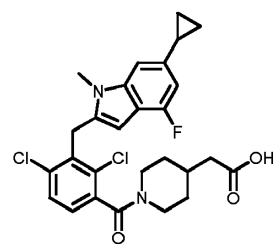
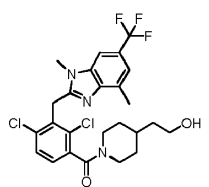
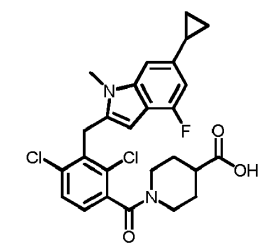
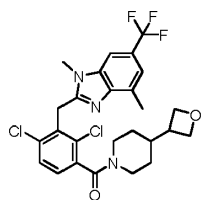
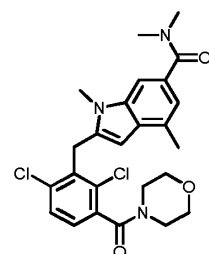
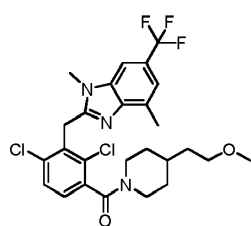
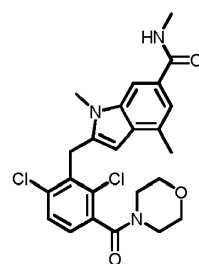
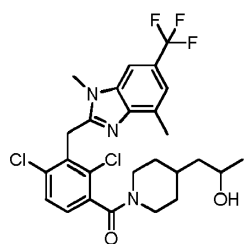
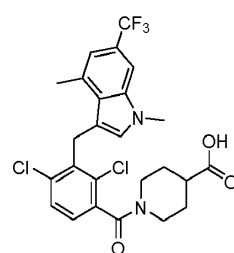
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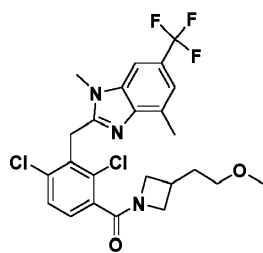
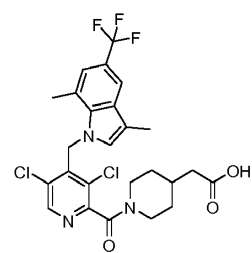
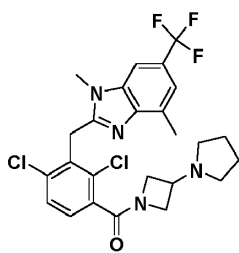
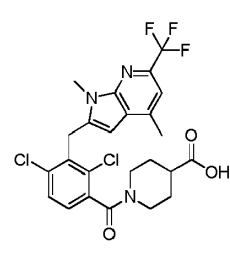
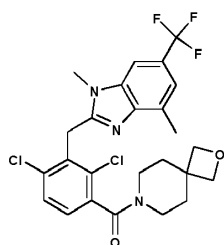
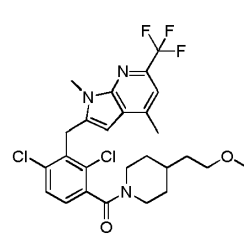
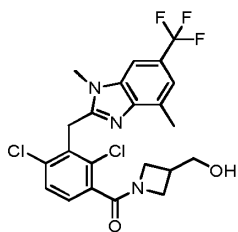
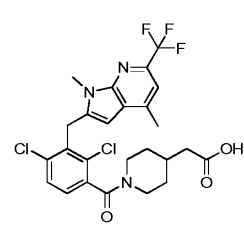
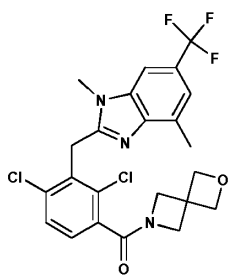
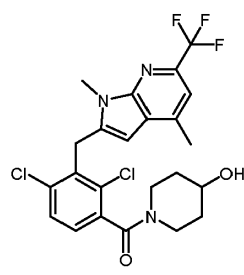
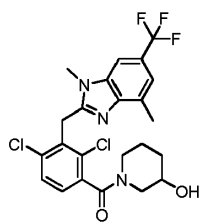
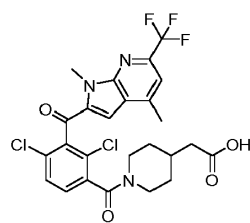
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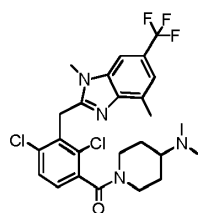
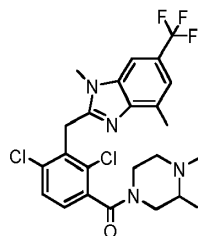
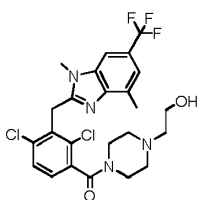
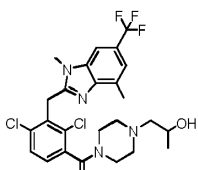
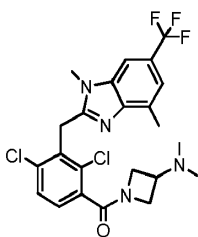
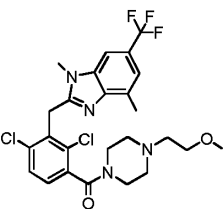
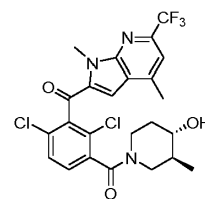
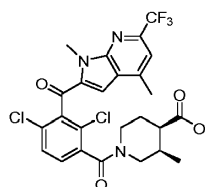
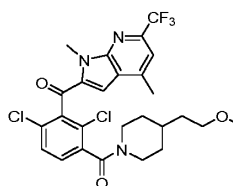
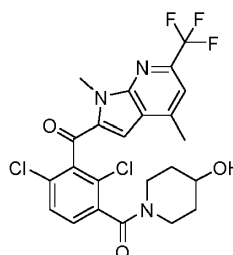
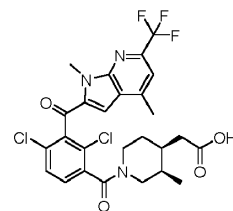
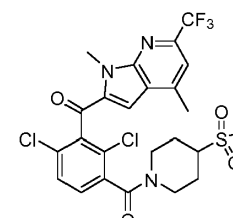
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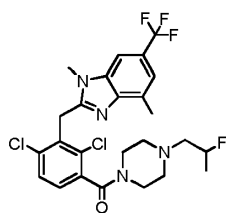
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**BJ-8****FG****BJ-9****FH****BJ-10****FI****BJ-11****FI-1****BJ-12****FJ****BJ-13****FK**

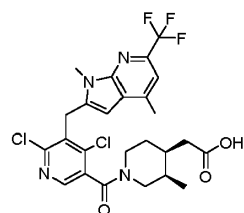
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**BJ-21****FR****BJ-23****FS****BJ-25****FT****BJ-27****FT-1****BJ-28****FT-3****BJ-29****FU**

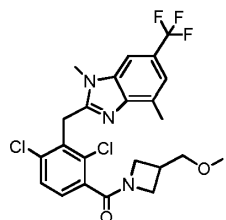
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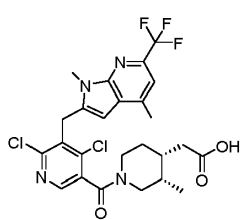
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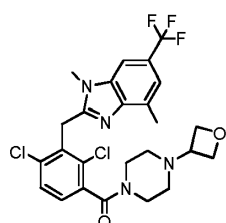
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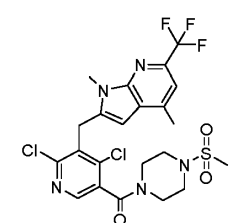
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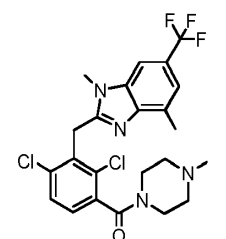
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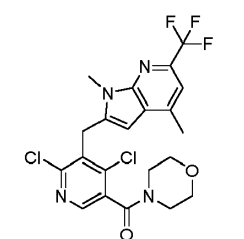
BJ-38



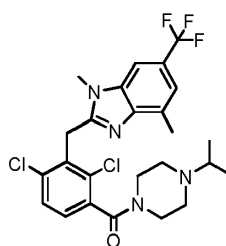
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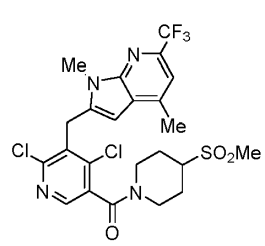
BJ-39



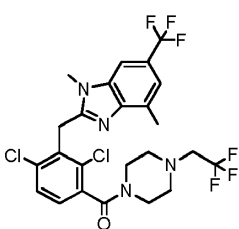
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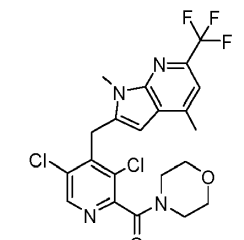
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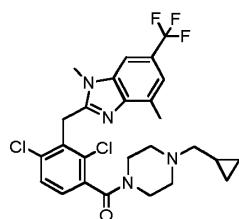
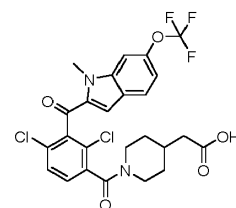
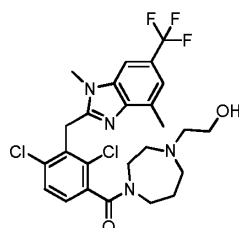
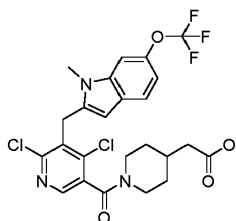
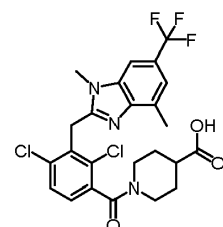
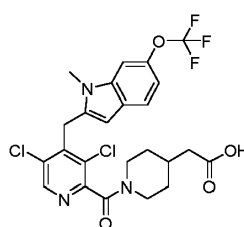
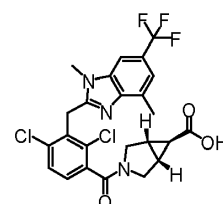
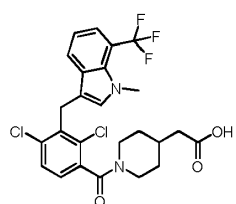
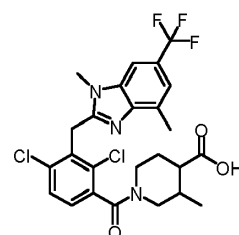
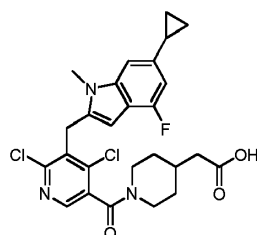
GG



BJ-41



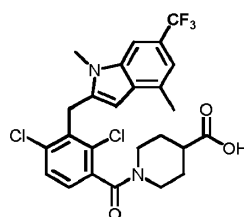
GH

**BJ-42****GI****BJ-43****GJ****BK****GK****BL****GL****BM****GN**

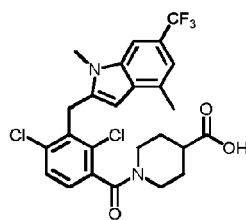
5

eller et farmasøytisk akseptabelt salt av samme.

13. Forbindelse ifølge krav 12, hvori forbindelsen er:

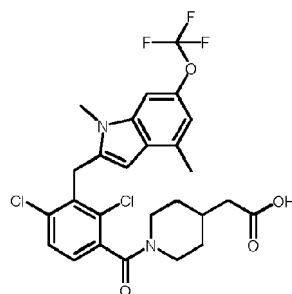
**B.**

14. Forbindelse ifølge krav 12, hvori forbindelsen er et farmasøytisk akseptabelt salt av:



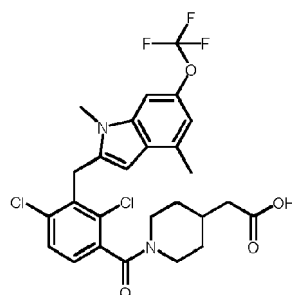
B.

15. Forbindelse ifølge krav 12, hvori forbindelsen er:



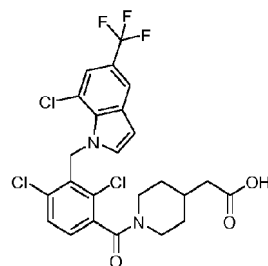
DP.

5 16. Forbindelse ifølge krav 12, hvori forbindelsen er et farmasøytisk akseptabelt salt av:



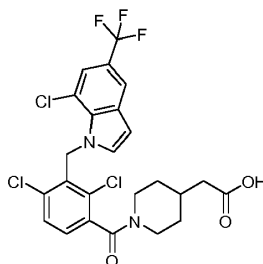
DP.

17. Forbindelse ifølge krav 12, hvori forbindelsen er:



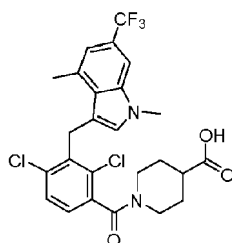
EI-5.

18. Forbindelse ifølge krav 12, hvori forbindelsen er et farmasøytisk akseptabelt salt av:



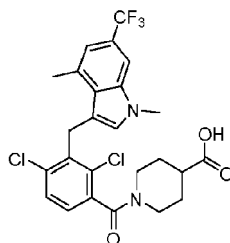
EI-5.

19. Forbindelse ifølge krav 12, hvori forbindelsen er:



FQ.

5 20. Forbindelse ifølge krav 12, hvori forbindelsen er et farmasøytisk akseptabelt salt av:



FQ.

21. Farmasøytisk blanding som omfatter:

- (i) en forbindelse ifølge ett av kravene 1 til 11, eller et farmasøytisk akseptabelt salt av samme, eller
 - 10 (ii) en forbindelse ifølge krav 12 eller et farmasøytisk akseptabelt salt av samme, eller
 - (iii) en forbindelse ifølge ett av kravene 13, 15, 17, eller 19, eller
 - (iv) et farmasøytisk akseptabelt salt av en forbindelse ifølge ett av kravene 14, 16, 18, eller 20,
- og ett eller flere farmasøytisk akseptable hjelpestoffer.